

Data Path : Z:\SVOASRV\HPCHEM1\BNA P\DATA\BP091019\  
 Data File : BP000188.D  
 Acq On : 11 Sep 2019 00:40  
 Operator : HP/JU  
 Sample : K4633-15  
 Misc :  
 ALS Vial : 24 Sample Multiplier: 1

**Instrument :**  
 BNA\_P  
**ClientSampled :**  
 F2D14

**Manual Integrations**  
**APPROVED**  
 mohammad  
 9/12/2019 10:26:00 AM

Quant Time: Sep 11 04:36:07 2019  
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA\_P\METHODS\SOM-EPA-BP090519MA.M  
 Quant Title : SVOA CALIBRATION  
 QLast Update : Wed Sep 11 02:03:45 2019  
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.89	152	454126	20.00	ng/ul	0.00
18) Naphthalene-d8	8.18	136	1714368	20.00	ng/ul	0.00
36) Acenaphthene-d10	9.95	164	852975	20.00	ng/ul	0.00
62) Phenanthrene-d10	11.45	188	1435205	20.00	ng/ul	0.00
78) Chrysene-d12	14.15	240	1068537	20.00	ng/ul	0.00
86) Perylene-d12	15.70	264	1022899	20.00	ng/ul	0.02

**System Monitoring Compounds**

3) 1,4-Dioxane-d8	2.61	96	57896	4.52	ng/uL	0.00
5) Phenol-d5	6.51	99	1019007	26.34	ng/ul	0.00
7) Bis-(2-Chloroethyl)ether-d	6.58	67	527112	26.06	ng/ul	0.00
9) 2-Chlorophenol-d4	6.66	132	837470	26.77	ng/ul	0.00
13) 4-Methylphenol-d8	7.26	113	730338	25.13	ng/ul	0.00
19) Nitrobenzene-d5	7.45	128	399770	30.71	ng/ul	0.00
22) 2-Nitrophenol-d4	7.78	143	437265	34.50	ng/ul	0.00
26) 2,4-Dichlorophenol-d3	8.02	165	730171	27.15	ng/ul	0.00
29) 4-Chloroaniline-d4	8.23	131	593709	17.85	ng/ul	0.00
44) Dimethylphthalate-d6	9.63	166	1948227	30.93	ng/ul	0.00
47) Acenaphthylene-d8	9.79	160	2269401	29.16	ng/ul	0.00
52) 4-Nitrophenol-d4	10.03	143	305112	27.11	ng/ul	0.00
58) Fluorene-d10	10.47	176	1544977	28.88	ng/ul	0.00
63) 4,6-Dinitro-2-methylphenol	10.53	200	98398	14.92	ng/ul	0.00
71) Anthracene-d10	11.50	188	2055622	29.58	ng/ul	0.00
79) Pyrene-d10	12.90	212	1972604	31.04	ng/ul	0.00
90) Benzo(a)pyrene-d12	15.61	264	1609646	30.36	ng/ul	0.02

**Target Compounds**

					Ovalue
6) Phenol	6.53	94	185757	4.608	ng/ul 99
45) Dimethylphthalate	9.66	163	631940	9.943	ng/ul 100
77) Fluoranthene	12.68	202	220880	2.580	ng/ul 98
80) Pyrene	12.92	202	191551	2.376	ng/ul 94
83) Benzo(a)anthracene	14.13	228	128133	1.657	ng/ul 94
85) Chrysene	14.17	228	211199m	2.976	ng/ul
88) Benzo(b)fluoranthene	15.24	252	388974m	5.737	ng/ul
89) Benzo(k)fluoranthene	15.26	252	96129m	1.504	ng/ul
91) Benzo(a)pyrene	15.63	252	180787	2.842	ng/ul# 64
92) Indeno(1,2,3-cd)pyrene	17.29	276	210327	2.874	ng/ul# 88
93) Dibenzo(a,h)anthracene	17.30	278	84453	1.393	ng/ul# 68
94) Benzo(g,h,i)perylene	17.76	276	238948	3.916	ng/ul# 88

(#) = qualifier out of range (m) = manual integration (+) = signals summed

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