

Data Path : Z:\SVOASRV\HPCHEM1\BNA_P\DATA\BP122420\
 Data File : BP004431.D
 Acq On : 24 Dec 2020 12:59
 Operator : CG/JU
 Sample : L5132-02
 Misc :
 ALS Vial : 5 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 A54R7

Quant Time: Dec 24 13:59:48 2020
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_P\METHODS\SOM-EPA-BP121820MA.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Thu Dec 24 13:54:26 2020
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.82	152	243152	20.00	ng/ul	0.00
18) Naphthalene-d8	10.62	136	899989	20.00	ng/ul	0.00
36) Acenaphthene-d10	14.47	164	519728	20.00	ng/ul	0.00
62) Phenanthrene-d10	17.23	188	1125439	20.00	ng/ul	0.00
78) Chrysene-d12	21.32	240	1313942	20.00	ng/ul	0.00
86) Perylene-d12	23.68	264	1422272	20.00	ng/ul	0.00

System Monitoring Compounds

3) 1,4-Dioxane-d8	3.32	96	30846	5.28	ng/uL	0.00
5) Phenol-d5	6.99	99	545667	25.85	ng/ul	0.00
7) Bis-(2-Chloroethyl)ether-d	7.15	67	310503	26.23	ng/ul	0.00
9) 2-Chlorophenol-d4	7.35	132	449074	26.01	ng/ul	0.00
13) 4-Methylphenol-d8	8.53	113	409446	24.35	ng/ul	0.00
19) Nitrobenzene-d5	8.97	128	216849	27.84	ng/ul	0.00
22) 2-Nitrophenol-d4	9.70	143	237059	27.91	ng/ul	0.00
26) 2,4-Dichlorophenol-d3	10.24	165	422817	26.84	ng/ul	0.00
29) 4-Chloroaniline-d4	10.75	131	534057	29.90	ng/ul	0.00
44) Dimethylphthalate-d6	13.87	166	1160282	27.89	ng/ul	0.00
47) Acenaphthylene-d8	14.16	160	1568135	30.35	ng/ul	0.00
52) 4-Nitrophenol-d4	14.67	143	171124	22.77	ng/ul	0.00
58) Fluorene-d10	15.47	176	1031303	28.34	ng/ul	0.00
63) 4,6-Dinitro-2-methylphenol	15.59	200	138094	19.02	ng/ul	0.00
71) Anthracene-d10	17.33	188	1636159	29.43	ng/ul	0.00
79) Pyrene-d10	19.57	212	1959629	27.92	ng/ul	0.00
90) Benzo(a)pyrene-d12	23.53	264	2352567	30.03	ng/ul	0.00

Target Compounds

					Ovalue
6) Phenol	7.02	94	108200	5.095 ng/ul	98
45) Dimethylphthalate	13.92	163	190820	4.578 ng/ul	98

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

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