

Data Path : Z:\SVOASRV\HPCHEM1\BNA P\DATA\BP090819\
 Data File : BP000120.D
 Acq On : 09 Sep 2019 04:00
 Operator : HP/JU
 Sample : K4664-01 2X
 Misc :
 ALS Vial : 18 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 OR-03-090619

Manual Integrations
 APPROVED

mohammad
 9/10/2019 2:51:24 PM

Quant Time: Sep 09 08:16:04 2019
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA P\METHODS\8270-BP083019.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Aug 30 17:47:17 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.89	152	459347	20.00	ng	0.00
21) Naphthalene-d8	8.18	136	1714530	20.00	ng	0.00
39) Acenaphthene-d10	9.93	164	929638	20.00	ng	-0.01
64) Phenanthrene-d10	11.43	188	1656033	20.00	ng	0.00
76) Chrysene-d12	14.10	240	1304653	20.00	ng	-0.01
87) Perylene-d12	15.63	264	1325898	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.51	112	1554297	53.96	ng	0.00
7) Phenol-d6	6.51	99	2035235	55.97	ng	0.00
23) Nitrobenzene-d5	7.45	82	1056523	37.81	ng	0.00
42) 2,4,6-Tribromophenol	10.73	330	473238	60.39	ng	0.00
45) 2-Fluorobiphenyl	9.25	172	2184870	39.25	ng	0.00
79) Terphenyl-d14	13.03	244	2481092	41.45	ng	0.00
Target Compounds						
49) Acenaphthylene	9.79	152	180489	2.233	ng	99
50) Dimethylphthalate	9.65	163	274019	4.325	ng	99
71) Phenanthrene	11.45	178	400833	4.856	ng	96
72) Anthracene	11.50	178	192758	2.364	ng	96
75) Fluoranthene	12.66	202	1050992	11.717	ng	92
78) Pyrene	12.89	202	1124927	11.721	ng	97
81) Benzo(a)anthracene	14.09	228	557954m	6.600	ng	
83) Chrysene	14.12	228	551604	6.898	ng	97
86) Indeno(1,2,3-cd)pyrene	17.17	276	298393	2.978	ng	94
88) Benzo(b)fluoranthene	15.17	252	766223m	9.461	ng	
89) Benzo(k)fluoranthene	15.20	252	183760m	2.479	ng	
90) Benzo(a)pyrene	15.56	252	528376	7.227	ng	98
92) Benzo(a,h,i)perylene	17.63	276	294129	4.065	ng	95

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

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