

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP120721\
 Data File : BP008258.D
 Acq On : 07 Dec 2021 19:32
 Operator : CG/JU
 Sample : M4920-04
 Misc :
 ALS Vial : 7 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 FDR-20211202-WC-W

Quant Time: Dec 08 01:10:13 2021
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\8270E-BP112521.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Dec 03 14:39:24 2021
 Response via : Initial Calibration

Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.781	152	62069	20.000	ng	-0.02
21) Naphthalene-d8	10.581	136	221290	20.000	ng	-0.02
39) Acenaphthene-d10	14.434	164	136604	20.000	ng	-0.01
64) Phenanthrene-d10	17.198	188	290657	20.000	ng	-0.01
76) Chrysene-d12	21.304	240	357964	20.000	ng	-0.01
86) Perylene-d12	23.692	264	426563	20.000	ng	-0.02
System Monitoring Compounds						
5) 2-Fluorophenol	5.375	112	558079	144.768	ng	-0.01
7) Phenol-d6	6.969	99	655659	125.992	ng	-0.02
23) Nitrobenzene-d5	8.946	82	525152	107.914	ng	-0.02
42) 2,4,6-Tribromophenol	15.934	330	285369	163.439	ng	-0.02
45) 2-Fluorobiphenyl	13.051	172	1037012	110.267	ng	-0.02
79) Terphenyl-d14	19.769	244	1785620	104.251	ng	-0.01
Target Compounds						
						Qvalue
80) Butylbenzylphthalate	20.445	149	25320	2.484	ng	95
81) Benzo(a)anthracene	21.286	228	55484	2.339	ng	98
83) Chrysene	21.339	228	65743	2.912	ng	98
84) Bis(2-ethylhexyl)phtha...	21.216	149	36462	2.356	ng	99
85) Di-n-octyl phthalate	22.133	149	80335	3.056	ng	# 99
88) Benzo(b)fluoranthene	22.969	252	87816	3.415	ng	97
89) Benzo(k)fluoranthene	23.016	252	92270	3.665	ng	95
90) Benzo(a)pyrene	23.592	252	73398	3.340	ng	96

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

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