

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP120321\
 Data File : BP008209.D
 Acq On : 03 Dec 2021 17:27
 Operator : CG/JU
 Sample : M4798-09
 Misc :
 ALS Vial : 8 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 DUP-111721

Manual Integrations
 APPROVED

Reviewed By :Jagrut Upadhyay 12/06/2021
 Supervised By :Sohil Jodhani 12/06/2021

Quant Time: Dec 03 18:17:17 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	7.793	152	103052	20.000	ng	0.00
21) Naphthalene-d8	10.593	136	369747	20.000	ng	# 0.00
39) Acenaphthene-d10	14.451	164	235709	20.000	ng	0.00
64) Phenanthrene-d10	17.216	188	501648	20.000	ng	0.00
76) Chrysene-d12	21.310	240	465286	20.000	ng	0.00
86) Perylene-d12	23.704	264	414986	20.000	ng	-0.01
System Monitoring Compounds						
5) 2-Fluorophenol	5.387	112	480106	75.012	ng	0.00
7) Phenol-d6	6.981	99	754732	87.353	ng	0.00
23) Nitrobenzene-d5	8.957	82	587025	72.195	ng	0.00
42) 2,4,6-Tribromophenol	15.957	330	415421	137.887	ng	0.00
45) 2-Fluorobiphenyl	13.069	172	1216763	74.982	ng	0.00
79) Terphenyl-d14	19.780	244	2039251	91.597	ng	0.00
Target Compounds						Qvalue
38) 1-Methylnaphthalene	12.487	142	242304	18.582	ng	# 97
52) Acenaphthene	14.516	154	74932	6.035	ng	# 90
55) Dibenzofuran	14.857	168	60027	2.829	ng	# 1
58) Fluorene	15.504	166	114185m	6.623	ng	
71) Phenanthrene	17.257	178	205476	7.758	ng	# 87
78) Pyrene	19.580	202	171880	3.478	ng	# 93

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

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