

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP033122\
 Data File : BP009644.D
 Acq On : 31 Mar 2022 17:50
 Operator : CG/JU
 Sample : N2198-01
 Misc :
 ALS Vial : 13 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 RO-VNJ-202

Quant Time: Apr 01 04:10:41 2022
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\8270E-BP031722.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu Mar 31 04:36:51 2022
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.740	152	302665	20.000	ng	0.00
21) Naphthalene-d8	10.534	136	1225505	20.000	ng	0.00
39) Acenaphthene-d10	14.387	164	761587	20.000	ng	0.00
64) Phenanthrene-d10	17.157	188	1509032	20.000	ng	0.00
76) Chrysene-d12	21.274	240	1410017	20.000	ng	0.00
86) Perylene-d12	23.663	264	1515874	20.000	ng	0.01
System Monitoring Compounds						
5) 2-Fluorophenol	5.346	112	2045810	118.013	ng	0.00
7) Phenol-d6	6.934	99	2622999	111.885	ng	0.00
23) Nitrobenzene-d5	8.899	82	1679700	72.835	ng	0.00
42) 2,4,6-Tribromophenol	15.892	330	1156709	92.040	ng	0.00
45) 2-Fluorobiphenyl	13.010	172	3524113	64.148	ng	0.00
79) Terphenyl-d14	19.739	244	4943295	62.926	ng	0.00
Target Compounds						
						Qvalue
10) Phenol	6.963	94	60764	2.584	ng	77
71) Phenanthrene	17.198	178	233010	2.885	ng	98
75) Fluoranthene	19.186	202	496935	5.081	ng	99
78) Pyrene	19.539	202	461931	4.972	ng	99
81) Benzo(a)anthracene	21.263	228	249302	2.691	ng	94
83) Chrysene	21.310	228	254108	2.897	ng	95
84) Bis(2-ethylhexyl)phtha...	21.192	149	153886	2.796	ng	# 99
88) Benzo(b)fluoranthene	22.933	252	335661	3.709	ng	# 95
90) Benzo(a)pyrene	23.557	252	234136	3.024	ng	# 94

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

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