

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP040122\
 Data File : BP009660.D
 Acq On : 01 Apr 2022 16:22
 Operator : CG/JU
 Sample : N2167-02
 Misc :
 ALS Vial : 10 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 RAIL-PILE-1

Quant Time: Apr 01 22:44:20 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	7.728	152	218528	20.000	ng	-0.01
21) Naphthalene-d8	10.522	136	853760	20.000	ng	-0.01
39) Acenaphthene-d10	14.381	164	524300	20.000	ng	-0.01
64) Phenanthrene-d10	17.145	188	1115734	20.000	ng	-0.01
76) Chrysene-d12	21.269	240	1180988	20.000	ng	0.00
86) Perylene-d12	23.645	264	1286025	20.000	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.340	112	2008232	160.448	ng	0.00
7) Phenol-d6	6.928	99	2254190	133.174	ng	0.00
23) Nitrobenzene-d5	8.887	82	1698818	105.739	ng	-0.01
42) 2,4,6-Tribromophenol	15.887	330	1174584	135.761	ng	0.00
45) 2-Fluorobiphenyl	12.999	172	3808607	100.703	ng	-0.01
79) Terphenyl-d14	19.728	244	6395933	97.207	ng	-0.01
Target Compounds						
						Qvalue
31) Naphthalene	10.569	128	211660	4.813	ng	98
52) Acenaphthene	14.446	154	224917	7.920	ng	97
55) Dibenzofuran	14.787	168	103397	2.167	ng	99
58) Fluorene	15.440	166	150412	4.009	ng	98
71) Phenanthrene	17.187	178	460832	7.716	ng	100
73) Carbazole	17.569	167	240738	4.174	ng	99
75) Fluoranthene	19.175	202	202195	2.796	ng	98

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

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