

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP021122\
 Data File : BP009120.D
 Acq On : 11 Feb 2022 19:04
 Operator : CG/JU
 Sample : N1426-03
 Misc :
 ALS Vial : 11 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 POST-EX-03-0207

Quant Time: Feb 11 23:33:45 2022
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\8270E-BP020722.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Feb 07 15:13:17 2022
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.793	152	206366	20.000	ng	-0.01
21) Naphthalene-d8	10.587	136	833209	20.000	ng	-0.02
39) Acenaphthene-d10	14.434	164	593064	20.000	ng	-0.01
64) Phenanthrene-d10	17.192	188	1306583	20.000	ng	-0.01
76) Chrysene-d12	21.292	240	1466948	20.000	ng	0.00
86) Perylene-d12	23.680	264	1433619	20.000	ng	-0.01
System Monitoring Compounds						
5) 2-Fluorophenol	5.387	112	1421756	122.895	ng	0.00
7) Phenol-d6	6.987	99	1829237	119.998	ng	0.00
23) Nitrobenzene-d5	8.957	82	1115333	75.489	ng	-0.01
42) 2,4,6-Tribromophenol	15.939	330	1032969	130.450	ng	0.00
45) 2-Fluorobiphenyl	13.057	172	3173799	74.916	ng	-0.01
79) Terphenyl-d14	19.757	244	5230528	64.121	ng	-0.01
Target Compounds						
						Qvalue
71) Phenanthrene	17.233	178	229806	3.248	ng	96
75) Fluoranthene	19.210	202	568571	7.171	ng	98
78) Pyrene	19.563	202	513733	4.923	ng	98
81) Benzo(a)anthracene	21.274	228	289176	3.059	ng	99
83) Chrysene	21.327	228	218090	2.393	ng	99
88) Benzo(b)fluoranthene	22.957	252	318807	3.612	ng	99
90) Benzo(a)pyrene	23.574	252	237360	3.226	ng	96

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

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