

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP033122\
 Data File : BP009646.D
 Acq On : 31 Mar 2022 18:58
 Operator : CG/JU
 Sample : N2199-03
 Misc :
 ALS Vial : 15 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 RO-ETGI-371

Quant Time: Apr 01 04:11:09 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	350704	20.000	ng	0.00
21) Naphthalene-d8	10.534	136	1390073	20.000	ng	0.00
39) Acenaphthene-d10	14.392	164	859033	20.000	ng	0.00
64) Phenanthrene-d10	17.157	188	1704141	20.000	ng	0.00
76) Chrysene-d12	21.274	240	1632633	20.000	ng	0.00
86) Perylene-d12	23.663	264	1823515	20.000	ng	0.01
System Monitoring Compounds						
5) 2-Fluorophenol	5.346	112	2175298	108.294	ng	0.00
7) Phenol-d6	6.934	99	2739606	100.852	ng	0.00
23) Nitrobenzene-d5	8.899	82	1757027	67.169	ng	0.00
42) 2,4,6-Tribromophenol	15.892	330	1136317	80.160	ng	0.00
45) 2-Fluorobiphenyl	13.010	172	3542443	57.167	ng	0.00
79) Terphenyl-d14	19.739	244	5212537	57.306	ng	0.00
Target Compounds						
						Qvalue
10) Phenol	6.963	94	67395	2.473	ng	# 73
71) Phenanthrene	17.198	178	269893	2.959	ng	99
75) Fluoranthene	19.186	202	620567	5.619	ng	98
78) Pyrene	19.539	202	580772	5.398	ng	99
81) Benzo(a)anthracene	21.257	228	299016	2.788	ng	97
83) Chrysene	21.310	228	307073	3.023	ng	98
88) Benzo(b)fluoranthene	22.939	252	415769	3.819	ng	# 94
90) Benzo(a)pyrene	23.557	252	264788	2.843	ng	# 94

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

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