

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN081924\
 Data File : BN033499.D
 Acq On : 20 Aug 2024 10:58
 Operator : MA/JU
 Sample : P3650-04
 Misc :
 ALS Vial : 30 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 RW7-SP303-20240814

Quant Time: Aug 20 11:38:06 2024
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_N\Methods\8270-SIM-BN081924.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Aug 19 23:32:18 2024
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.552	152	8308	0.400	ng	0.00
7) Naphthalene-d8	10.314	136	21989	0.400	ng	# 0.00
13) Acenaphthene-d10	14.189	164	10862	0.400	ng	0.00
19) Phenanthrene-d10	16.942	188	22983	0.400	ng	0.00
29) Chrysene-d12	21.148	240	13365	0.400	ng	0.00
35) Perylene-d12	23.317	264	12335	0.400	ng	0.00
System Monitoring Compounds						
4) 2-Fluorophenol	5.183	112	2411	0.091	ng	0.00
5) Phenol-d6	6.736	99	1748	0.056	ng	0.00
8) Nitrobenzene-d5	8.691	82	5071	0.278	ng	0.00
11) 2-Methylnaphthalene-d10	11.911	152	8954	0.285	ng	0.00
14) 2,4,6-Tribromophenol	15.688	330	1259	0.216	ng	0.00
15) 2-Fluorobiphenyl	12.809	172	13483	0.304	ng	0.00
27) Fluoranthene-d10	18.979	212	18508	0.335	ng	0.00
31) Terphenyl-d14	19.593	244	13457	0.443	ng	0.00
Target Compounds						Qvalue
2) 1,4-Dioxane	3.190	88	459	0.048	ng	# 1
32) Benzo(a)anthracene	21.130	228	1365	0.028	ng	# 89
33) Chrysene	21.184	228	1668	0.035	ng	# 92
34) Bis(2-ethylhexyl)phtha...	21.094	149	3000	0.098	ng	99
36) Indeno(1,2,3-cd)pyrene	25.478	276	1742	0.034	ng	99
37) Benzo(b)fluoranthene	22.674	252	1663	0.036	ng	# 25
38) Benzo(k)fluoranthene	22.715	252	1608	0.035	ng	# 26
39) Benzo(a)pyrene	23.224	252	1172	0.031	ng	# 1
40) Dibenzo(a,h)anthracene	25.490	278	1364	0.033	ng	# 47
41) Benzo(g,h,i)perylene	26.127	276	1450	0.033	ng	# 66

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

