

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN092922\
 Data File : BN021956.D
 Acq On : 29 Sep 2022 15:54
 Operator : CG/JU
 Sample : N4846-01
 Misc :
 ALS Vial : 6 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 TT15811-20220923

Quant Time: Sep 30 02:45:22 2022
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_N\METHODS\8270-SIM-BN092922.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Sep 30 02:27:15 2022
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	8.033	152	4011	0.400	ng	0.00
7) Naphthalene-d8	10.866	136	12162	0.400	ng	# 0.01
13) Acenaphthene-d10	14.686	164	6468	0.400	ng	0.00
19) Phenanthrene-d10	17.437	188	13023	0.400	ng	0.00
29) Chrysene-d12	21.627	240	9938	0.400	ng	# 0.00
35) Perylene-d12	24.131	264	8021	0.400	ng	0.00
System Monitoring Compounds						
4) 2-Fluorophenol	5.563	112	1255	0.118	ng	0.00
5) Phenol-d6	7.188	99	894	0.066	ng	0.00
8) Nitrobenzene-d5	9.211	82	2652	0.261	ng	0.00
11) 2-Methylnaphthalene-d10	12.452	152	8406	0.320	ng	0.00
14) 2,4,6-Tribromophenol	16.184	330	818	0.295	ng	0.00
15) 2-Fluorobiphenyl	13.317	172	8136	0.284	ng	0.00
27) Fluoranthene-d10	19.471	212	12946	0.367	ng	0.00
31) Terphenyl-d14	20.060	244	9310	0.424	ng	0.00
Target Compounds						
34) Bis(2-ethylhexyl)phtha...	21.529	149	1847	0.080	ng	Qvalue 98
36) Indeno(1,2,3-cd)pyrene	26.766	276	930	0.022	ng	# 63
41) Benzo(g,h,i)perylene	27.572	276	1379	0.037	ng	# 74

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

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