

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN042221\
 Data File : BN014367.D
 Acq On : 21 Apr 2021 23:26
 Operator : CG/JU
 Sample : M2122-02
 Misc :
 ALS Vial : 7 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 BP-VPB-RW8-GW-58-60

Quant Time: Apr 22 01:55:16 2021
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_N\METHODS\8270-SIM-BN042221.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Apr 21 13:57:02 2021
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.692	152	6900	0.40	ng	0.00
7) Naphthalene-d8	10.467	136	24363	0.40	ng	0.00
13) Acenaphthene-d10	14.321	164	12280	0.40	ng	0.00
19) Phenanthrene-d10	17.067	188	25945	0.40	ng	0.00
29) Chrysene-d12	21.261	240	23682	0.40	ng	# 0.00
36) Perylene-d12	23.486	264	19926	0.40	ng	0.00
System Monitoring Compounds						
4) 2-Fluorophenol	5.292	112	2602	0.18	ng	0.00
5) Phenol-d6	6.863	99	2411	0.14	ng	0.00
8) Nitrobenzene-d5	8.832	82	5913	0.29	ng	0.00
11) 2-Methylnaphthalene-d10	12.057	152	9387	0.26	ng	0.00
14) 2,4,6-Tribromophenol	15.818	330	1276	0.32	ng	0.00
15) 2-Fluorobiphenyl	12.938	172	15058	0.31	ng	0.00
27) Fluoranthene-d10	19.104	212	22930	0.33	ng	0.00
31) Terphenyl-d14	19.705	244	30613	0.61	ng	0.00
Target Compounds						
25) Phenanthrene	17.104	178	1531	0.02	ng	# 95
28) Fluoranthene	19.134	202	1739	0.02	ng	# 93
30) Pyrene	19.496	202	1662	0.02	ng	98
32) Benzo(a)anthracene	21.250	228	1950	0.03	ng	# 93
33) Chrysene	21.292	228	2224	0.03	ng	93
34) Bis(2-ethylhexyl)phtha...	21.186	149	57101	2.58	ng	99
35) Indeno(1,2,3-cd)pyrene	25.724	276	1800	0.02	ng	98
37) Benzo(b)fluoranthene	22.818	252	1821	0.02	ng	# 23
38) Benzo(k)fluoranthene	22.863	252	1694	0.02	ng	# 19
39) Benzo(a)pyrene	23.386	252	1313	0.02	ng	# 1
40) Dibenzo(a,h)anthracene	25.735	278	1605	0.02	ng	# 42
41) Benzo(g,h,i)perylene	26.402	276	1807	0.02	ng	# 76

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

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