

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN042621\
 Data File : BN014442.D
 Acq On : 26 Apr 2021 13:46
 Operator : CG/JU
 Sample : M2164-03
 Misc :
 ALS Vial : 6 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 BP-VPB-RW8-GW-258-260

Quant Time: Apr 26 14:16:14 2021
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_N\METHODS\8270-SIM-BN042221.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Apr 26 12:30:05 2021
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.684	152	8439	0.40	ng	0.00
7) Naphthalene-d8	10.454	136	31355	0.40	ng	0.00
13) Acenaphthene-d10	14.308	164	16572	0.40	ng	0.00
19) Phenanthrene-d10	17.055	188	37144	0.40	ng	# 0.00
29) Chrysene-d12	21.261	240	34039	0.40	ng	# 0.00
36) Perylene-d12	23.482	264	26593	0.40	ng	0.00
System Monitoring Compounds						
4) 2-Fluorophenol	5.284	112	1324	0.08	ng	0.00
5) Phenol-d6	6.855	99	1064	0.05	ng	0.00
8) Nitrobenzene-d5	8.819	82	3042	0.12	ng	0.00
11) 2-Methylnaphthalene-d10	12.049	152	9318	0.20	ng	0.00
14) 2,4,6-Tribromophenol	15.805	330	639	0.12	ng	0.00
15) 2-Fluorobiphenyl	12.938	172	14675	0.22	ng	0.01
27) Fluoranthene-d10	19.094	212	22146	0.23	ng	0.00
31) Terphenyl-d14	19.700	244	23625	0.33	ng	0.00
Target Compounds						
2) 1,4-Dioxane	3.247	88	3083	0.30	ng	# 73
32) Benzo(a)anthracene	21.239	228	2210	0.02	ng	93
33) Chrysene	21.292	228	2449	0.02	ng	94
34) Bis(2-ethylhexyl)phtha...	21.186	149	34559	1.09	ng	100

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

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