

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN090421\
 Data File : BN016198.D
 Acq On : 04 Sep 2021 18:04
 Operator : CG/JU
 Sample : M3604-01
 Misc :
 ALS Vial : 13 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 PW-B6-L41-083021

Manual Integrations
 APPROVED

mohammad
 9/7/2021 11:40:13 AM

Quant Time: Sep 06 06:19:47 2021
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_N\METHODS\8270-SIM-BN090421.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Sep 06 02:08:58 2021
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.873	152	13241	0.400	ng	0.00
7) Naphthalene-d8	10.660	136	71817	0.400	ng	0.00
13) Acenaphthene-d10	14.497	164	44340	0.400	ng	0.00
19) Phenanthrene-d10	17.249	188	86477	0.400	ng	# 0.00
29) Chrysene-d12	21.429	240	93586	0.400	ng	# 0.00
35) Perylene-d12	23.823	264	98220	0.400	ng	0.00
System Monitoring Compounds						
4) 2-Fluorophenol	5.466	112	3851	0.103	ng	0.00
5) Phenol-d6	7.043	99	2886	0.060	ng	0.00
8) Nitrobenzene-d5	9.025	82	14538	0.252	ng	0.00
11) 2-Methylnaphthalene-d10	12.252	152	29543	0.258	ng	0.00
14) 2,4,6-Tribromophenol	15.994	330	5935	0.300	ng	0.00
15) 2-Fluorobiphenyl	13.127	172	42398	0.254	ng	0.00
27) Fluoranthene-d10	19.276	212	80733	0.311	ng	0.00
31) Terphenyl-d14	19.872	244	79811	0.337	ng	0.00
Target Compounds						Qvalue
2) 1,4-Dioxane	3.373	88	992m	0.234	ng	
9) Naphthalene	10.711	128	10925	0.057	ng	# 95
12) 2-Methylnaphthalene	12.327	142	55062	0.429	ng	# 99
17) Acenaphthene	14.560	154	10225	0.081	ng	# 38
18) Fluorene	15.550	166	46562	0.294	ng	# 72
24) Pentachlorophenol	16.908	266	376	0.382	ng	# 75
25) Phenanthrene	17.285	178	329366	1.360	ng	# 92
28) Fluoranthene	19.306	202	24006	0.084	ng	# 85
30) Pyrene	19.668	202	57340	0.188	ng	# 79
34) Bis(2-ethylhexyl)phtha...	21.344	149	206709	0.977	ng	98

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

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