

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN090421\
 Data File : BN016215.D
 Acq On : 05 Sep 2021 05:33
 Operator : CG/JU
 Sample : M3620-03
 Misc :
 ALS Vial : 28 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 BP-TT-VPB-RW7-GW-640-642

Manual Integrations
 APPROVED

mohammad
 9/7/2021 11:40:33 AM

Quant Time: Sep 06 06:22:01 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	7.864	152	14538	0.400	ng	0.00
7) Naphthalene-d8	10.660	136	78294	0.400	ng	# 0.00
13) Acenaphthene-d10	14.497	164	47347	0.400	ng	0.00
19) Phenanthrene-d10	17.236	188	94377	0.400	ng	#-0.01
29) Chrysene-d12	21.429	240	99530	0.400	ng	# 0.00
35) Perylene-d12	23.813	264	101387	0.400	ng	-0.01
System Monitoring Compounds						
4) 2-Fluorophenol	5.457	112	3014	0.073	ng	0.00
5) Phenol-d6	7.034	99	2489	0.047	ng	0.00
8) Nitrobenzene-d5	9.025	82	17020	0.271	ng	0.00
11) 2-Methylnaphthalene-d10	12.247	152	33241	0.267	ng	0.00
14) 2,4,6-Tribromophenol	15.981	330	4163	0.197	ng	-0.01
15) 2-Fluorobiphenyl	13.114	172	48058	0.269	ng	-0.01
27) Fluoranthene-d10	19.271	212	97610	0.344	ng	-0.01
31) Terphenyl-d14	19.867	244	92161	0.366	ng	0.00
Target Compounds						Qvalue
2) 1,4-Dioxane	3.373	88	8958m	1.926	ng	
9) Naphthalene	10.711	128	11318	0.055	ng	97
12) 2-Methylnaphthalene	12.323	142	11803	0.084	ng	99
17) Acenaphthene	14.560	154	6280	0.047	ng	# 66
18) Fluorene	15.537	166	17014	0.101	ng	# 80
24) Pentachlorophenol	16.896	266	1716	0.429	ng	98
25) Phenanthrene	17.285	178	96813	0.366	ng	# 92
28) Fluoranthene	19.301	202	8325	0.027	ng	# 91
30) Pyrene	19.668	202	16504	0.051	ng	# 80
34) Bis(2-ethylhexyl)phtha...	21.333	149	55034	0.245	ng	98

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

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