

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN072724\
 Data File : BN033042.D
 Acq On : 27 Jul 2024 20:51
 Operator : MA/JU
 Sample : P3375-04
 Misc :
 ALS Vial : 58 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 EB-01-072524

Manual Integrations
 APPROVED

Reviewed By :Jagrut Upadhyay 07/29/2024
 Supervised By :mohammad ahmed 07/30/2024

Quant Time: Jul 29 01:08:09 2024
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_N\Methods\8270-SIM-BN071124.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Sun Jul 21 23:33:21 2024
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.582	152	1537m	0.400	ng	0.05
7) Naphthalene-d8	10.372	136	5925m	0.400	ng	0.07
13) Acenaphthene-d10	14.190	164	4065	0.400	ng	0.01
19) Phenanthrene-d10	16.970	188	9060	0.400	ng	# 0.02
29) Chrysene-d12	21.178	240	8529	0.400	ng	0.01
35) Perylene-d12	23.361	264	9022	0.400	ng	0.00
System Monitoring Compounds						
4) 2-Fluorophenol	5.242	112	838m	0.216	ng	0.07
5) Phenol-d6	7.091	99	698m	0.145	ng	0.31
8) Nitrobenzene-d5	8.984	82	1625m	0.364	ng	0.25
11) 2-Methylnaphthalene-d10	11.954	152	2982	0.329	ng	0.04
14) 2,4,6-Tribromophenol	15.753	330	518	0.298	ng	0.03
15) 2-Fluorobiphenyl	12.822	172	4413	0.292	ng	0.01
27) Fluoranthene-d10	18.997	212	9966	0.425	ng	0.00
31) Terphenyl-d14	19.587	244	8586	0.491	ng	0.00
Target Compounds						
2) 1,4-Dioxane	3.097	88	311	0.163	ng	# 24
33) Chrysene	21.214	228	1256m	0.039	ng	
41) Benzo(g,h,i)perylene	26.308	276	886	0.028	ng	# 32

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

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