

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN090921\
 Data File : BN016297.D
 Acq On : 10 Sep 2021 19:10
 Operator : CG/JU
 Sample : M3620-07
 Misc :
 ALS Vial : 45 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 BP-TT-VPB-RW7-GW-680-682

Manual Integrations
 APPROVED

mohammad
 9/13/2021 10:21:40 AM

Quant Time: Sep 11 02:30:43 2021
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_N\METHODS\8270-SIM-BN090921.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Sep 10 14:34:37 2021
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.864	152	11554	0.400	ng	0.00
7) Naphthalene-d8	10.660	136	59725	0.400	ng	0.00
13) Acenaphthene-d10	14.484	164	34476	0.400	ng	-0.01
19) Phenanthrene-d10	17.237	188	71379	0.400	ng	0.00
29) Chrysene-d12	21.429	240	63360	0.400	ng	0.00
35) Perylene-d12	23.813	264	38971	0.400	ng	0.00
System Monitoring Compounds						
4) 2-Fluorophenol	5.466	112	2549	0.086	ng	0.00
5) Phenol-d6	7.052	99	2063	0.056	ng	0.02
8) Nitrobenzene-d5	9.013	82	12811	0.267	ng	0.00
11) 2-Methylnaphthalene-d10	12.247	152	24356	0.261	ng	0.00
14) 2,4,6-Tribromophenol	15.982	330	3872	0.274	ng	0.00
15) 2-Fluorobiphenyl	13.114	172	35762	0.265	ng	0.00
27) Fluoranthene-d10	19.271	212	65917	0.314	ng	0.00
31) Terphenyl-d14	19.867	244	65753	0.414	ng	0.00
Target Compounds						
2) 1,4-Dioxane	3.373	88	9999m	2.439	ng	Qvalue
24) Pentachlorophenol	16.896	266	107	0.127	ng	# 72
34) Bis(2-ethylhexyl)phtha...	21.334	149	32594	0.268	ng	100

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

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