

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN092021\
 Data File : BN016352.D
 Acq On : 20 Sep 2021 19:59
 Operator : CG/JU
 Sample : M3729-13
 Misc :
 ALS Vial : 17 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 BP-TT-VPB-RW7-GW-820-822

Manual Integrations
 APPROVED

mohammad
 9/21/2021 11:30:33 AM

Quant Time: Sep 21 01:34:16 2021
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_N\METHODS\8270-SIM-BN091721.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Sep 20 11:09:23 2021
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.689	152	25492	0.400	ng	0.00
7) Naphthalene-d8	10.457	136	115691	0.400	ng	0.00
13) Acenaphthene-d10	14.307	164	64799	0.400	ng	0.00
19) Phenanthrene-d10	17.054	188	130353	0.400	ng	0.00
29) Chrysene-d12	21.259	240	140724	0.400	ng	0.00
35) Perylene-d12	23.522	264	137282	0.400	ng	0.00
System Monitoring Compounds						
4) 2-Fluorophenol	5.319	112	7344	0.115	ng	0.00
5) Phenol-d6	6.868	99	5321	0.069	ng	0.00
8) Nitrobenzene-d5	8.822	82	17575	0.256	ng	0.00
11) 2-Methylnaphthalene-d10	12.047	152	40404	0.222	ng	0.00
14) 2,4,6-Tribromophenol	15.804	330	6848	0.360	ng	0.00
15) 2-Fluorobiphenyl	12.937	172	56125	0.221	ng	0.00
27) Fluoranthene-d10	19.098	212	105769	0.280	ng	0.00
31) Terphenyl-d14	19.708	244	102718	0.315	ng	0.00
Target Compounds						Qvalue
24) Pentachlorophenol	16.713	266	552	0.023	ng	98
25) Phenanthrene	17.103	178	7418m	0.020	ng	
34) Bis(2-ethylhexyl)phtha...	21.196	149	177038	0.726	ng	99

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

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