

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN100421\
 Data File : BN016732.D
 Acq On : 05 Oct 2021 05:49
 Operator : CG/JU
 Sample : M3829-08DL 2X
 Misc :
 ALS Vial : 25 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 RE125D2-20210917DL

Manual Integrations
 APPROVED

mohammad
 10/5/2021 12:08:40 PM

Quant Time: Oct 05 06:27:59 2021
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_N\METHODS\8270-SIM-BN100421.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Oct 04 17:52:33 2021
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.642	152	28433	0.40	ng	0.00
7) Naphthalene-d8	10.406	136	128121	0.40	ng	0.00
13) Acenaphthene-d10	14.268	164	67225	0.40	ng	0.00
19) Phenanthrene-d10	17.029	188	129038	0.40	ng	0.00
29) Chrysene-d12	21.238	240	117668	0.40	ng	# 0.00
35) Perylene-d12	23.488	264	121395	0.40	ng	0.00
System Monitoring Compounds						
4) 2-Fluorophenol	5.263	112	3134	0.04	ng	0.00
5) Phenol-d6	6.821	99	1794	0.02	ng	0.00
8) Nitrobenzene-d5	8.784	82	8683	0.10	ng	0.00
11) 2-Methylnaphthalene-d10	12.007	152	23941	0.13	ng	0.00
14) 2,4,6-Tribromophenol	15.778	330	3335	0.20	ng	0.00
15) 2-Fluorobiphenyl	12.898	172	32466	0.12	ng	0.00
27) Fluoranthene-d10	19.068	212	56835	0.16	ng	0.00
31) Terphenyl-d14	19.678	244	47853	0.18	ng	0.00
Target Compounds						
2) 1,4-Dioxane	3.253	88	41331	3.11	ng	# 69
25) Phenanthrene	17.066	178	10886m	0.03	ng	
34) Bis(2-ethylhexyl)phtha...	21.174	149	8366	0.04	ng	99

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

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