

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM040221\
 Data File : BM029336.D
 Acq On : 03 Apr 2021 02:16
 Operator : CG/JU
 Sample : M1883-04
 Misc :
 ALS Vial : 24 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 V-211

Manual Integrations
 APPROVED

mohammad
 4/5/2021 11:55:47 AM

Quant Time: Apr 03 02:57:46 2021
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\8270-BM040121.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu Apr 01 16:52:48 2021
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.698	152	102638	20.00	ng	0.00
21) Naphthalene-d8	10.475	136	389087	20.00	ng	-0.01
39) Acenaphthene-d10	14.333	164	217282	20.00	ng	0.00
64) Phenanthrene-d10	17.068	188	436154	20.00	ng	0.00
76) Chrysene-d12	21.245	240	443188	20.00	ng	0.00
86) Perylene-d12	23.450	264	470747	20.00	ng	-0.01
System Monitoring Compounds						
5) 2-Fluorophenol	5.299	112	880471	141.64	ng	0.00
7) Phenol-d6	6.875	99	1051324	119.80	ng	0.00
23) Nitrobenzene-d5	8.851	82	827434	97.50	ng	0.00
42) 2,4,6-Tribromophenol	15.821	330	338595	155.91	ng	0.00
45) 2-Fluorobiphenyl	12.951	172	1586862	101.75	ng	-0.01
79) Terphenyl-d14	19.698	244	2217837	95.05	ng	0.00
Target Compounds						
						Qvalue
31) Naphthalene	10.528	128	246145	12.00	ng	99
32) Benzoic acid	9.739	122	29532	7.16	ng	92
37) 2-Methylnaphthalene	12.139	142	108266	7.49	ng	98
38) 1-Methylnaphthalene	12.363	142	69625m	5.09	ng	
70) Pentachlorophenol	16.727	266	289	5.77	ng	93

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

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