

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM041921\
 Data File : BM029552.D
 Acq On : 19 Apr 2021 22:12
 Operator : CG/JU
 Sample : M2074-13
 Misc :
 ALS Vial : 19 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 VNPH2-B-9

Manual Integrations
 APPROVED

mohammad
 4/21/2021 8:35:44 AM

Quant Time: Apr 20 04:17:09 2021
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\8270-BM040121.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Apr 16 11:56:50 2021
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.663	152	110094	20.00	ng	-0.01
21) Naphthalene-d8	10.445	136	426612	20.00	ng	-0.01
39) Acenaphthene-d10	14.304	164	214073	20.00	ng	-0.01
64) Phenanthrene-d10	17.051	188	350203	20.00	ng	0.00
76) Chrysene-d12	21.227	240	293438	20.00	ng	0.00
86) Perylene-d12	23.432	264	324792	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.269	112	754185	113.11	ng	0.00
7) Phenol-d6	6.845	99	960085	101.99	ng	0.00
23) Nitrobenzene-d5	8.822	82	693261	74.50	ng	0.00
42) 2,4,6-Tribromophenol	15.798	330	203138	94.94	ng	0.00
45) 2-Fluorobiphenyl	12.927	172	1156282	75.25	ng	0.00
79) Terphenyl-d14	19.680	244	991592	64.18	ng	0.00
Target Compounds						Qvalue
10) Phenol	6.875	94	24865	2.69	ng	85
50) Dimethylphthalate	13.768	163	42332	2.74	ng	98
75) Fluoranthene	19.103	202	72608	3.22	ng	100
78) Pyrene	19.468	202	63868	3.13	ng	96
88) Benzo(b)fluoranthene	22.768	252	49325m	2.31	ng	

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

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