

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP062121\
 Data File : BP005996.D
 Acq On : 21 Jun 2021 16:16
 Operator : CG/JU
 Sample : M2709-18
 Misc :
 ALS Vial : 10 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 SPLP-AMND-COMP-O(CL-24)

Manual Integrations
 APPROVED

mohammad
 6/22/2021 10:15:33 AM

Quant Time: Jun 21 17:01:48 2021
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\8270E-BP060821.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Jun 16 11:35:30 2021
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.922	152	65156	20.00	ng	0.00
21) Naphthalene-d8	10.728	136	239767	20.00	ng	0.00
39) Acenaphthene-d10	14.551	164	148158	20.00	ng	0.00
64) Phenanthrene-d10	17.298	188	316613	20.00	ng	0.00
76) Chrysene-d12	21.369	240	382805	20.00	ng	-0.01
86) Perylene-d12	23.780	264	435477	20.00	ng	-0.01
System Monitoring Compounds						
5) 2-Fluorophenol	5.499	112	342116	84.25	ng	0.00
7) Phenol-d6	7.099	99	252555	47.99	ng	0.00
23) Nitrobenzene-d5	9.093	82	445958	91.19	ng	0.00
42) 2,4,6-Tribromophenol	16.040	330	396399	155.91	ng	-0.01
45) 2-Fluorobiphenyl	13.181	172	1001951	88.80	ng	0.00
79) Terphenyl-d14	19.845	244	1977224	96.72	ng	-0.01
Target Compounds						
32) Benzoic acid	10.105	122	1116m	6.68	ng	Qvalue

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

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