

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP061621\
 Data File : BP005950.D
 Acq On : 16 Jun 2021 14:46
 Operator : CG/JU
 Sample : M2678-06RX
 Misc :
 ALS Vial : 9 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 0239-AMND-COMP-D(CL-01A)RX

Manual Integrations
 APPROVED

mohammad
 6/17/2021 12:16:23 PM

Quant Time: Jun 16 15:20:32 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	7.928	152	67979	20.000	ng	0.00
21) Naphthalene-d8	10.734	136	255063	20.000	ng	0.00
39) Acenaphthene-d10	14.557	164	162286	20.000	ng	0.00
64) Phenanthrene-d10	17.304	188	356685	20.000	ng	0.00
76) Chrysene-d12	21.374	240	438026	20.000	ng	0.00
86) Perylene-d12	23.786	264	503102	20.000	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.505	112	12486	2.947	ng	0.00
7) Phenol-d6	7.105	99	236669	43.101	ng	0.00
23) Nitrobenzene-d5	9.093	82	303654	58.367	ng	0.00
42) 2,4,6-Tribromophenol	16.057	330	657	0.236	ng	0.00
45) 2-Fluorobiphenyl	13.187	172	662911	53.639	ng	0.00
79) Terphenyl-d14	19.851	244	1151879	49.241	ng	0.00
Target Compounds						
						Qvalue
75) Fluoranthene	19.304	202	69563	2.674	ng	99
78) Pyrene	19.657	202	83427	2.728	ng	99
81) Benzo(a)anthracene	21.363	228	78431	2.480	ng	97
83) Chrysene	21.410	228	71993m	2.351	ng	
88) Benzo(b)fluoranthene	23.051	252	97898m	2.841	ng	
90) Benzo(a)pyrene	23.680	252	76144	2.350	ng	# 95

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

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0239-AMND-COMP-D(CL-01A)RX

Manual Integrations
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