

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP092421\
 Data File : BP007163.D
 Acq On : 24 Sep 2021 17:41
 Operator : CG/JU
 Sample : M3875-15
 Misc :
 ALS Vial : 4 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 20882

Manual Integrations
 APPROVED

mohammad
 9/27/2021 3:04:18 PM

Quant Time: Sep 24 18:10:23 2021
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\8270E-BP092121.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu Sep 23 11:17:25 2021
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.887	152	251640	20.000	ng	0.00
21) Naphthalene-d8	10.687	136	987967	20.000	ng	# 0.00
39) Acenaphthene-d10	14.522	164	521488	20.000	ng	0.00
64) Phenanthrene-d10	17.269	188	1198778	20.000	ng	# 0.00
76) Chrysene-d12	21.351	240	1141689	20.000	ng	# 0.00
86) Perylene-d12	23.768	264	203452	20.000	ng	# 0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.469	112	1920806	127.775	ng	0.00
7) Phenol-d6	7.063	99	1979149	96.327	ng	0.00
23) Nitrobenzene-d5	9.051	82	1480681	87.280	ng	0.00
42) 2,4,6-Tribromophenol	16.022	330	1032152	152.666	ng	0.00
45) 2-Fluorobiphenyl	13.145	172	3417211	93.877	ng	0.00
79) Terphenyl-d14	19.827	244	4602939	77.258	ng	0.00
Target Compounds						
						Qvalue
2) 1,4-Dioxane	3.364	88	390537m	64.242	ng	
10) Phenol	7.093	94	1420705	71.721	ng	90
20) 3+4-Methylphenols	8.657	107	853853	46.243	ng	85
32) Benzoic acid	10.046	122	340538	33.365	ng	93
84) Bis(2-ethylhexyl)phtha...	21.263	149	211300	4.899	ng	99
88) Benzo(b)fluoranthene	23.027	252	45357	3.730	ng	# 16

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

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