

Data Path : Z:\SVOASRV\HPCHEM1\BNA P\DATA\BP111519\
 Data File : BP001052.D
 Acq On : 15 Nov 2019 21:26
 Operator : JU
 Sample : K5832-11
 Misc :
 ALS Vial : 15 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 6-(6-8)

Manual Integrations
 APPROVED

mohammad
 11/18/2019 3:06:33 PM

Quant Time: Nov 16 04:39:54 2019
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA P\METHODS\8270-BP110619.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Nov 06 17:07:13 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	8.37	152	213842	20.00	ng	0.00
21) Naphthalene-d8	10.39	136	784836	20.00	ng	-0.01
39) Acenaphthene-d10	13.25	164	456119	20.00	ng	-0.02
64) Phenanthrene-d10	15.65	188	923557	20.00	ng	-0.01
76) Chrysene-d12	19.28	240	789897	20.00	ng	-0.02
87) Perylene-d12	21.06	264	866863	20.00	ng	-0.02
System Monitoring Compounds						
5) 2-Fluorophenol	6.31	112	1133886	96.51	ng	0.01
7) Phenol-d6	7.82	99	1459517	98.24	ng	0.00
23) Nitrobenzene-d5	9.24	82	921844	64.97	ng	-0.02
42) 2,4,6-Tribromophenol	14.55	330	623257	114.46	ng	-0.01
45) 2-Fluorobiphenyl	12.17	172	2064462	68.17	ng	-0.02
79) Terphenyl-d14	17.93	244	2765423	69.64	ng	-0.01
Target Compounds						
10) Phenol	7.83	94	45745	2.815	ng	90
50) Dimethylphthalate	12.83	163	345442	10.242	ng	98
71) Phenanthrene	15.68	178	114853	2.430	ng	99
75) Fluoranthene	17.39	202	206452	3.864	ng	97
78) Pyrene	17.69	202	244584	4.381	ng	100
83) Chrysene	19.31	228	94769	2.029	ng	98
88) Benzo(b)fluoranthene	20.57	252	106777m	2.066	ng	

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

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