

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP071822\
 Data File : BP011015.D
 Acq On : 19 Jul 2022 04:27
 Operator : CG/JU
 Sample : PB146234BL
 Misc :
 ALS Vial : 31 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 SBLK234

Quant Time: Jul 19 05:07:54 2022
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP071422.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Tue Jul 19 03:25:54 2022
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.675	152	421437	20.000	ng/u1	0.00
20) Naphthalene-d8	10.469	136	1645721	20.000	ng/u1	# 0.00
38) Acenaphthene-d10	14.339	164	862924	20.000	ng/u1	0.00
64) Phenanthrene-d10	17.104	188	1675648	20.000	ng/u1	# 0.00
79) Chrysene-d12	21.227	240	1681644	20.000	ng/u1	# 0.00
88) Perylene-d12	23.580	264	1775127	20.000	ng/u1	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.181	96	68500	5.860	ng/uL	0.00
4) Pyridine-d5	3.587	84	793859	26.273	ng/u1	0.00
7) Phenol-d5	6.857	99	985762	28.495	ng/u1	0.00
9) Bis-(2-Chloroethyl)eth...	7.022	67	663583	29.830	ng/u1	0.00
11) 2-Chlorophenol-d4	7.210	132	839513	30.033	ng/u1	0.00
15) 4-Methylphenol-d8	8.393	113	759057	27.408	ng/u1	0.00
21) Nitrobenzene-d5	8.840	128	374617	31.236	ng/u1	0.00
24) 2-Nitrophenol-d4	9.557	143	358949	30.247	ng/u1	0.00
28) 2,4-Dichlorophenol-d3	10.092	165	613208	28.255	ng/u1	0.00
31) 4-Chloroaniline-d4	10.616	131	1043004	29.152	ng/u1	0.00
46) Dimethylphthalate-d6	13.757	166	1852823	31.416	ng/u1	0.00
49) Acenaphthylene-d8	14.028	160	2220535	31.910	ng/u1	0.00
54) 4-Nitrophenol-d4	14.557	143	266905	22.361	ng/u1	0.00
60) Fluorene-d10	15.339	176	1556822	31.055	ng/u1	0.00
65) 4,6-Dinitro-2-methylph...	15.463	200	187194	22.576	ng/u1	0.00
73) Anthracene-d10	17.204	188	2385334	33.485	ng/u1	0.00
81) Pyrene-d10	19.463	212	2699195	30.562	ng/u1	0.00
92) Benzo(a)pyrene-d12	23.427	264	2809778	32.956	ng/u1	0.00

Target Compounds Qvalue

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

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