

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP071023\
 Data File : BP016179.D
 Acq On : 11 Jul 2023 04:50
 Operator : MA/JU
 Sample : PB153948BL
 Misc :
 ALS Vial : 3 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 SBLK948

Manual Integrations
 APPROVED

Reviewed By :Yogesh Patel 07/11/2023
 Supervised By :mohammad ahmed 07/11/2023

Quant Time: Jul 11 05:39:02 2023
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP062623.MA.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Mon Jul 10 23:19:51 2023
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	8.011	152	291029	20.000	ng/u1	0.00
20) Naphthalene-d8	10.840	136	1336840	20.000	ng/u1	# 0.00
38) Acenaphthene-d10	14.663	164	867535	20.000	ng/u1	0.00
64) Phenanthrene-d10	17.428	188	1914854	20.000	ng/u1	# 0.00
79) Chrysene-d12	21.545	240	1579130	20.000	ng/u1	# 0.00
88) Perylene-d12	24.074	264	1475824	20.000	ng/u1	-0.02
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.340	96	48595	6.008	ng/uL	0.00
4) Pyridine-d5	3.787	84	587624	28.818	ng/u1	0.00
7) Phenol-d5	7.205	99	761757	30.592	ng/u1	0.00
9) Bis-(2-Chloroethyl)eth...	7.334	67	541773m	35.167	ng/u1	-0.01
11) 2-Chlorophenol-d4	7.540	132	617440	32.848	ng/u1	0.00
15) 4-Methylphenol-d8	8.758	113	659151	33.508	ng/u1	0.00
21) Nitrobenzene-d5	9.210	128	311108	30.451	ng/u1	0.00
24) 2-Nitrophenol-d4	9.946	143	337313	28.288	ng/u1	0.00
28) 2,4-Dichlorophenol-d3	10.528	165	498371	23.454	ng/u1	0.02
31) 4-Chloroaniline-d4	11.034	131	776059	28.431	ng/u1	0.01
46) Dimethylphthalate-d6	14.081	166	2143907	31.973	ng/u1	0.00
49) Acenaphthylene-d8	14.357	160	2373313	31.486	ng/u1	0.00
54) 4-Nitrophenol-d4	15.022	143	187647m	21.206	ng/u1	0.03
60) Fluorene-d10	15.675	176	1803502	32.665	ng/u1	0.00
65) 4,6-Dinitro-2-methylph...	15.863	200	217835	18.498	ng/u1	0.00
73) Anthracene-d10	17.539	188	2748688	31.425	ng/u1	0.00
81) Pyrene-d10	19.769	212	3194647	30.982	ng/u1	-0.01
92) Benzo(a)pyrene-d12	23.898	264	2320927	31.175	ng/u1	-0.02

Target Compounds Qvalue

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

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